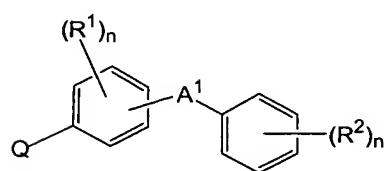


**Amendments to the Claims:**

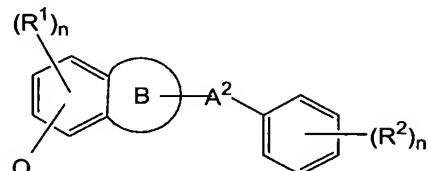
This listing of claims will replace all prior versions, and listings, of claims in the application:

**Listing of Claims:**

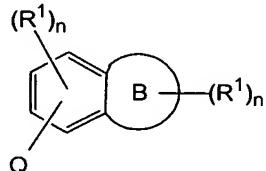
1. (Original) A compound having the formula:



(I)



(II)



(III)

or a pharmaceutically acceptable salt or solvate thereof, wherein:

Q is  $\text{CF}_3\text{SO}_2$ ,  $\text{CF}_3\text{SO}_2\text{NR}^3$ ,  $\text{CF}_3\text{SO}_2\text{R}^4$  or  $\text{CF}_3\text{SO}_2\text{N}(\text{R}^3)\text{R}^4$ , wherein  $\text{R}^3$  is H, alkoxy, acyl or  $\text{C}_1\text{-C}_3$  alkyl, each of which may be substituted or unsubstituted, and  $\text{R}^4$  is methylene which may be substituted or unsubstituted;

each  $\text{R}^1$  is independently  $\text{C}_1\text{-C}_3$  alkyl,  $\text{C}_1\text{-C}_3$  haloalkyl, CN,  $(\text{C}=\text{O})\text{OR}$ ,  $(\text{C}=\text{O})\text{R}^5$ , H, halo,  $\text{O}(\text{C}=\text{O})\text{R}$ , OR, OH, NHR,  $\text{NH}(\text{C}=\text{O})\text{OR}$ ,  $\text{NH}(\text{C}=\text{O})\text{R}^5$ ,  $\text{NO}_2$ ,  $\text{NHSO}_2\text{R}^5$ ,  $\text{SO}_2\text{R}^5$ ,  $\text{R}^4\text{SO}_2\text{CF}_3$  or tetrazole, wherein  $\text{R}^5$  is  $\text{CF}_3$ ,  $\text{C}_1\text{-C}_3$  alkyl, NHR and wherein R is H,  $\text{C}_1\text{-C}_3$  alkyl, aryl or heteroaryl, which may be substituted or unsubstituted;

each  $\text{R}^2$  is independently  $\text{C}_1\text{-C}_3$  alkyl,  $\text{C}_1\text{-C}_3$  haloalkyl, CN,  $(\text{C}=\text{O})\text{OR}$ ,  $(\text{C}=\text{O})\text{R}^5$ , H, halo,  $\text{O}(\text{C}=\text{O})\text{R}$ , OR, OH, NHR,  $\text{NH}(\text{C}=\text{O})\text{OR}$ ,  $\text{NH}(\text{C}=\text{O})\text{R}^5$ ,  $\text{NO}_2$ ,  $\text{NHSO}_2\text{R}^5$ ,  $\text{SO}_2\text{R}^5$ , tetrazole, or  $\text{X}^1\text{-R}^6\text{-X}^2$ , wherein  $\text{X}^1$  is present or absent and if present is O, N,  $(\text{C}=\text{O})$ ,  $(\text{C}=\text{O})\text{NH}$ ,  $\text{NH}(\text{C}=\text{O})$ ,  $\text{SO}_2\text{NH}$ ,  $\text{NHSO}_2$ ,  $\text{R}^6$  is  $\text{C}_1\text{-C}_3$  alkylene which may be substituted or unsubstituted and  $\text{X}^2$  is  $\text{CF}_3$ ,  $(\text{C}=\text{O})\text{OR}$ ,  $(\text{C}=\text{O})\text{R}^5$ , H,  $\text{NH}(\text{C}=\text{O})\text{R}^5$ ,  $\text{NH}(\text{C}=\text{O})\text{OR}$ ,  $\text{NHSO}_2\text{R}^5$ ,  $\text{NRR}^3$ ,  $\text{O}(\text{C}=\text{O})\text{R}$ , OR,  $\text{SO}_2\text{R}^5$ , tetrazole;

each n is independently from 0 to 3;

ring B is an aryl, carbocyclic, heteroaryl, heterocyclic or phenyl ring which may be substituted or unsubstituted;

A<sup>1</sup> is a linkage in which the shortest path is 2-8 atoms in length wherein the atoms in the linkage are carbon which may be substituted or unsubstituted or the carbon replaced with a single nitrogen or oxygen, or combination of nitrogen, oxygen and sulfur provided no two heteroatoms are adjacently linked in a linear linkage; the linkage may be or may contain an aryl, carbocyclic, heteroaryl, heterocyclic or a phenyl ring, which may be directly in the linkage or appended to the linkage; the linkage may be acylalkyl, alkenylene, alkoxy, alkoxyalkyl, alkoxyamino, alkoxyarylalkoxy, alkoxyarylalkyl, alkoxyarylamino, alkoxyaryloxyalkyl, alkylamino, alkylaminoalkyl, alkylaminoarylaminooalkyl, alkylaryl, alkylarylalkyl, alkylarylarnino, alkylaryloxy, alkylene, alkylenediamine, alkylenedioxy, alkyloxy, alkyloxyaryl, alkyloxyarylalkyloxy, alkyloxyaryloxyalkyl, C<sub>1</sub>-C<sub>6</sub> alkylsulfonylamino, alkylthio, alkylthioalkyl, alkynylene, C<sub>1</sub>-C<sub>6</sub> N-sulfonamido, C<sub>3</sub>-C<sub>7</sub> N-amido, aminoalkyl, aminoalkylamino, aminoalkylarylalkyl, aminoalkylarylalkylamino, aminoalkylaryloxy, aminoalkyloxy, aminoaryl, aminoarylalkyl, aminoarylcarbonyl, aminoaryloxy, aminoaryloxyalkyl, aminoarylsulfonyl, aryl, arylamino, ortho or para aryldioxy, substituted meta-aryldioxy, aryldiamine, aryloxy, aryloxyalkyl, aryloxyamino, aryloxyaminoalkyl, aryloxycarbonyl, aryloxysulfonyl, benzimidazole, benzo[b]furan, benzo[b]thiophene, C<sub>3</sub>-C<sub>7</sub> C-amido, carbonylarylarnino, carbonylarylcarbonyl, carbonylaryloxy, chromene, cycloalkylene, disulfide, furan, haloalkyl, imidazole, imidazolidine, imidazoline, indole, isothiazole, isoxazole, morpholine, oxadiazole, oxazole, oxirane, parathiazine, phenothiazine, piperazine, piperidine, purine, pyran, pyrazine, pyrazole, pyrazolidine, pyrimidine, pyridine, pyrrole, pyrrolidine, quinoline, C<sub>2</sub>-C<sub>6</sub> S-sulfonamido, sulfonylalkyl, sulfonylarylarnino, sulfonylaryloxy, sulfonylarylsulfonyl, thiadiazole, thiazole, thiophene, triazine, triazole, unsubstituted azeridine, C<sub>3</sub>-C<sub>6</sub> ureido, which may be substituted or unsubstituted; and

A<sup>2</sup> is a linkage in which the shortest path is 0-6 atoms in length wherein the atoms in the linkage are carbon which may be substituted or unsubstituted or the carbon replaced with a single nitrogen, oxygen or sulfur, or combination of nitrogen, oxygen and sulfur; the linkage may be or may contain an aryl, carbocyclic, heteroaryl, heterocyclic or a phenyl ring, which

may be directly in the linkage or appended to the linkage; the linkage may be single atom C, O, S or N which may be substituted or unsubstituted; the linkage may be acylalkyl, alkenylene, alkoxy, alkoxyalkyl, alkoxyamino, alkoxyarylalkoxy, alkoxyarylalkyl, alkoxyarylamino, alkoxyaryloxyalkyl, alkylamino, alkylaminoalkyl, alkylaminoarylalkyl, alkylaryl, alkylarylalkyl, alkylarylamino, alkylaryloxy, alkylene, alkylenediamine, alkylenedioxy, alkyloxy, alkyloxyaryl, alkyloxyarylalkyloxy, alkyloxyaryloxyalkyl, alkylsulfonylamino, alkylthio, alkylthioalkyl, alkynylene, N-sulfonamido, N-amido, aminoalkyl, aminoalkylamino, aminoalkylarylalkyl, aminoalkylarylalkylamino, aminoalkylaryloxy, aminoalkyloxy, aminoaryl, aminoarylalkyl, aminoarylcarbonyl, aminoaryloxy, aminoaryloxyalkyl, aminoarylsulfonyl, aryl, arylamino, ortho or para aryldioxy, substituted meta-aryldioxy, aryldiamine, aryloxy, aryloxyalkyl, aryloxyamino, aryloxyaminoalkyl, aryloxycarbonyl, aryloxysulfonyl, benzimidazole, benzo[b]furan, benzo[b]thiophene, C-amido, carbonylarylamin, carbonylarylcarbonyl, carbonylaryloxy, chromene, cycloalkylene, furan, haloalkyl, imidazole, imidazolidine, imidazoline, indole, isothiazole, isoxazole, morpholine, oxadiazole, oxazole, oxirane, parathiazine, phenothiazine, piperazine, piperidine, purine, pyran, pyrazine, pyrazole, pyrazolidine, pyrimidine, pyridine, pyrrole, pyrrolidine, quinoline, sulfonamido, sulfonylalkyl, sulfonylarylamin, sulfonylaryloxy, sulfonylarylsulfonyl, thiadiazole, thiazole, thiophene, triazine, triazole, unsubstituted azeridine, ureido, which may be substituted or unsubstituted.

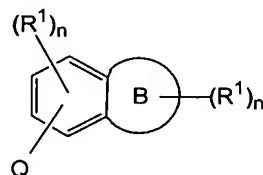
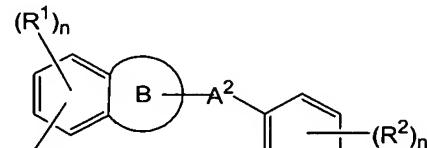
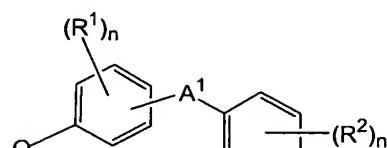
2. (Original) The compound of claim 1 wherein Q is  $\text{CF}_3\text{SO}_2$ .
3. (Original) The compound of claim 1 wherein Q is  $\text{CF}_3\text{SO}_2\text{NR}^3$ .
4. (Original) The compound of claim 1 wherein Q is  $\text{CF}_3\text{SO}_2\text{R}^4$ .
5. (Original) The compound of claim 1 wherein Q is  $\text{CF}_3\text{SO}_2\text{N}(\text{R}^3)\text{R}^4$ .
6. (Original) The compound of claim 1 wherein the compound has formula I.

7. (Original) The compound of claim 1 wherein the compound has formula II.

8. (Original) The compound of claim 1 wherein the compound has formula III.

9. (Original) The compound of claim 1 wherein R<sup>2</sup> is SO<sub>2</sub>R<sup>5</sup>, NHSO<sub>2</sub>R<sup>5</sup> or CF<sub>3</sub>SO<sub>2</sub>R<sup>4</sup>.

10. (Original) A compound having the formula:



or a pharmaceutically acceptable salt or solvate thereof, wherein:

Q is CF<sub>3</sub>SO<sub>2</sub>;

each R1 is independently CF<sub>3</sub>, (C=O)OR, (C=O)R5, H, halo, NHR, NH(C=O)OR, NH(C=O)R5, NHSO<sub>2</sub>R5, NO<sub>2</sub>, O(C=O)R, OH, OR, SO<sub>2</sub>R5 or tetrazole, wherein R5 is CF<sub>3</sub>, C1-C3 alkyl, NHR and wherein R is H, C1-C3 alkyl, aryl or heteroaryl, which may be substituted or unsubstituted;

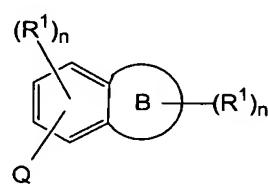
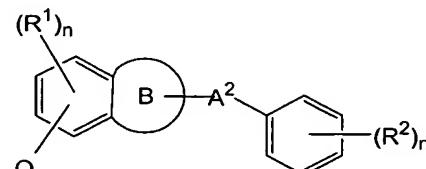
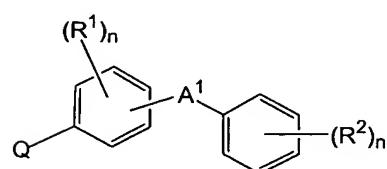
each R2 is independently (C=O)OR, (C=O)R5, NH(C=O)OR, NH(C=O)R5, NHR, NHSO<sub>2</sub>R5, NO<sub>2</sub>, -R6-(C=O)OR, -R6-NRR3, -R6-tetrazole, or tetrazole and R6 is C1-3 alkylene which may be substituted or unsubstituted;

each n is independently from 0 to 2;

ring B is phenyl or heteroaryl which may be substituted or unsubstituted; and

linkage A1 is C2-C4 alkoxy, C2-C4 alkoxyalkyl, C2-C4 alkylenedioxy, C2-C4 alkylaminoalkyl, C2-C4 alkylenediamine, C3-C4 C-amido, C3-C4 N-amido, C3-C4 ureido, C1-C3 N-sulfonamido, C2-C3 S-sulfonamido, aryldioxy, aryldiamine, aryl, alkylarylalkyl, imidazole, oxazole, oxadiazole, pyrazole, pyrazolidine, pyrrole or triazole.

11. (Original) A compound having the formula:



or a pharmaceutically acceptable salt or solvate thereof, wherein:

Q is  $\text{CF}_3\text{SO}_2\text{NR}^3$ ,  $\text{CF}_3\text{SO}_2\text{R}^4$  or  $\text{CF}_3\text{SO}_2\text{N}(\text{R}^3)\text{R}^4$ , wherein R3 is H, alkoxy, acyl or C1-C3 alkyl each of which may be substituted or unsubstituted and R4 is methylene which may be substituted or unsubstituted;

each R1 is independently  $\text{CF}_3$ ,  $(\text{C}=\text{O})\text{OR}$ ,  $(\text{C}=\text{O})\text{R}5$ , H, halo, NHR,  $\text{NH}(\text{C}=\text{O})\text{OR}$ ,  $\text{NH}(\text{C}=\text{O})\text{R}5$ ,  $\text{NHSO}_2\text{R}5$ ,  $\text{NO}_2$ ,  $\text{O}(\text{C}=\text{O})\text{R}$ , OH, OR,  $\text{SO}_2\text{R}5$  or tetrazole, wherein R5 is  $\text{CF}_3$ , C1-C3 alkyl, NHR and wherein R is H, C1-C3 alkyl, aryl or heteroaryl, which may be substituted or unsubstituted;

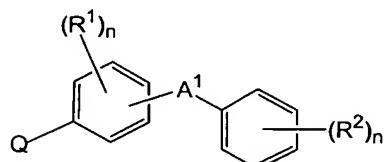
each R2 is independently  $(\text{C}=\text{O})\text{OR}$ ,  $(\text{C}=\text{O})\text{R}5$ ,  $\text{NH}(\text{C}=\text{O})\text{OR}$ ,  $\text{NH}(\text{C}=\text{O})\text{R}5$ , NHR,  $\text{NHSO}_2\text{R}5$ ,  $\text{NO}_2$ ,  $\text{SO}_2\text{R}5$ ,  $-\text{R}6-(\text{C}=\text{O})\text{OR}$ ,  $-\text{R}6-\text{NRR}3$ ,  $-\text{R}6$ -tetrazole, or tetrazole and R6 is C1-3 alkylene which may be substituted or unsubstituted;

each n is independently from 0 to 2;

ring B is phenyl or heteroaryl which may be substituted or unsubstituted; and

linkage A1 is C2-C4 alkoxy, C2-C4 alkoxyalkyl, C2-C4 alkylenedioxy, C2-C4 alkylaminoalkyl, C2-C4 alkylenediamine, C3-C4 C-amido, C3-C4 N-amido, C3-C4 ureido, C1-C3 N-sulfonamido, C2-C3 S-sulfonamido, aryldioxy, aryldiamine, aryl, alkylarylalkyl, imidazole, oxazole, oxadiazole, pyrazole, pyrazolidine, pyrrole or triazole.

12. (Original) A compound having the formula:



(I)

or a pharmaceutically acceptable salt or solvate thereof, wherein:

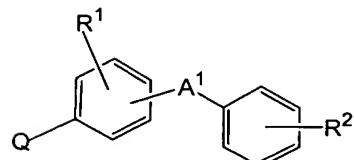
Q is  $\text{CF}_3\text{SO}_2$ ;

each R1 is independently H, NHR,  $\text{NO}_2$  or OR;

each R2 is independently  $(\text{C}=\text{O})\text{OR}$ ,  $\text{NHSO}_2\text{R}_5$ , or  $\text{SO}_2\text{R}_5$ ;

each n is independently from 0 to 2; and the linkage A1 is alkylarylalkyl, C2-C4 alkoxyalkyl, C2-C4 alkylenedioxy, aryl, aryldiamine, aryldioxy, or oxadiazole which may be substituted or unsubstituted or A1 is unsubstituted or monosubstituted C2-C4 N-amido.

13. (Original) The compound of claim 11 having the formula:



(IV)

or a pharmaceutically acceptable salt or solvate thereof, wherein:

R1 is H or  $\text{NO}_2$ ;

R2 is  $(\text{C}=\text{O})\text{OR}$ ,  $\text{NHSO}_2\text{R}_5$  or  $\text{SO}_2\text{R}_5$ ; and

the linkage A1 is C2-C4 alkoxyalkyl, aryldioxy, aryl, alkylarylalkyl or oxadiazole.

14. (Original) The compound of claim 12 wherein the compound is:  
Bis(4-Trifluoromethylsulfonylbenzyl) ether;  
4-Trifluoromethylsulfonylbenzyl 4-trifluoromethylsulfonylphenyl ether;  
N,N-Bis(4-trifluoromethylsulfonylbenzyl)benzamide;  
1,2-Bis(4-trifluoromethylsulfonylphenyl)ethane;  
N-(4-Trifluoromethylsulfonylbenzyl)-4-trifluoromethylsulfonylbenzamide;  
N-(4-Trifluoromethylsulfonylbenzyl)benzamide;  
3,5-Bis-(4-trifluoromethanesulfonyl-phenoxy)-benzoic acid methyl ester;  
[3,5-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-phenyl]-acetic acid methyl ester;  
3,5-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-benzoic acid methyl ester;  
1,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-cyclopentane;  
4-Methyl-2,6-bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-benzoic acid methyl ester;  
4-[2-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-ethoxy]-benzoic acid methyl ester;  
4-[3-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-phenoxy]-benzoic acid;  
1-(3,5-Bis-trifluoromethyl-phenyl)-5-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-1*H*-pyrazole-3-carboxylic acid methyl ester;  
{4-[4-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-benzenesulfonyl]-phenoxy}-acetic acid ethyl ester;  
4-[3-(4-Trifluoromethanesulfonyl-phenoxy)-phenoxy]-benzoic acid;  
{4-[4-(4-Trifluoromethanesulfonyl-phenoxy)-benzenesulfonyl]-phenoxy}-acetic acid ethyl ester;  
*N*-(3-Trifluoromethanesulfonyl-phenyl)-2-{2-[(3-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;  
*N*-(3-Trifluoromethanesulfonyl-phenyl)-2-{3-[(3-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;  
*N*-(3-Trifluoromethanesulfonyl-phenyl)-2-{4-[(3-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;

3,6-Bis-(morpholin-4-ylmethyl)-2,5-bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-benzene;

[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-dimethyl-amine;

*N*-(2-Ethylamino-5-trifluoromethanesulfonyl-phenyl)-2-(4-methanesulfonyl-phenyl)-acetamide;

2-Hydroxy-5-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-terephthalic acid diethyl ester;

{2-[(3-Trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetic acid;

{3-[(3-Trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetic acid;

{4-[(3-Trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetic acid;

3,5-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-benzamide;

3,5-Bis-(4-trifluoromethanesulfonyl-phenoxy)-benzoic acid;

*N*-(4-Trifluoromethanesulfonyl-phenyl)-2-{2-[(4-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;

*N*-(4-Trifluoromethanesulfonyl-phenyl)-2-{3-[(4-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;

*N*-(4-Trifluoromethanesulfonyl-phenyl)-2-{4-[(4-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-benzoic acid methyl ester;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-benzoic acid;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-*N*-pyridin-4-yl-benzamide;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-*N*-(4-methoxy-phenyl)-benzamide;

3-[4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-benzoylamino]-benzoic acid ethyl ester;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-*N*-(2-pyrrolidin-1-yl-ethyl)-benzamide;

*N*-Ethyl-4-(1-ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-benzamide;

1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazole-2-carboxylic acid;

[2-(Benzoyl-butyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid methyl ester;

*N*-Benzyl-*N*-[butylcarbamoyl-(4-trifluoromethanesulfonyl-phenyl)-methyl]-benzamide;

*N*-[Butylcarbamoyl-(4-trifluoromethanesulfonyl-phenyl)-methyl]-*N*-(2-hydroxyethyl)-benzamide;

[2-(Acetyl-cyclopropyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

[2-(Acetyl-methyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

[2-(Benzoyl-cyclohexyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

*N*-Cyclohexyl-*N*-[(2,6-dimethyl-phenylcarbamoyl)-(4-trifluoromethanesulfonyl-phenyl)-methyl]-benzamide;

[2-(Acetyl-propyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

[2-(Acetyl-cyclohexyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

{4-[4-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-benzenesulfonyl]-phenoxy}-acetic acid;

4-[2-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-ethoxy]-benzoic acid;

2,5-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-terephthalic acid diethyl ester;

1-[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-piperidine;

4-[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-morpholine;

[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-(2-nitro-phenyl)-amine;

1-(2-Nitro-phenylamino)-3-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propan-2-ol;

[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-(4-nitro-phenyl)-amine;

1-(4-Nitro-phenylamino)-3-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propan-2-ol;

4-[2-Hydroxy-3-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propylamino]-benzenesulfonamide; or

4-[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propylamino]-benzenesulfonamide.

15. (Original) The compound of claim 13 wherein the compound is:

1,2-Bis(4-trifluoromethylsulfonamidophenyl)ethane;

1,2-Bis(2-methyl-4-trifluoromethylsulfonamidophenyl)ethane;

1,3-Bis(4-trifluoromethylsulfonamidophenoxy)-2,2-dimethylpropane;

1,3-Bis(4-trifluoromethylsulfonamidophenoxy)propane;

1,4-Bis(4-trifluoromethylsulfonamidophenoxy)butane;

1,4-Bis(4-trifluoromethylsulfonamidophenoxy)benzene;

1-(4-Aminophenoxy)-4-trifluoromethylsulfonamidophenoxy benzene;

Bis(4-trifluoromethylsulfonamidophenyl) ether;

1,3-Bis(4-trifluoromethylsulfonamidophenoxy)benzene;

2,5-Bis(4-trifluoromethylsulfonamidophenyl)-(1,3,4)oxadiazole;

Bis(4-trifluoromethylsulfonamidophenyl)-1,4-diisopropylbenzene;

5-Trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid;

1-Methyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid;

(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetic acid;

1-Methyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid phenylamide;

5-Trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid phenylamide;

3-[(1-Methyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carbonyl)-amino]-benzoic acid;

3-[(5-Trifluoromethanesulfonylamino-1*H*-indole-2-carbonyl)-amino]-benzoic acid;

4-[(5-Trifluoromethanesulfonylamino-1*H*-indole-2-carbonyl)-amino]-benzoic acid;

4-[2-(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetylamino]-benzoic acid;

3-[2-(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetylamino]-benzoic acid;

4- {[2-(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetylamino]-methyl}-benzoic acid;

(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetic acid *tert*-butyl ester;

1-Methyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid ethyl ester;

6-Trifluoromethanesulfonylamino-naphthalene-2-carboxylic acid;

*N,N*-Bis[(6-carboxyl-naphthalen-2-yl)methyl] trifluoromethanesulfonamide;

6-[(Methyl-trifluoromethanesulfonyl-amino)-methyl]-naphthalene-2-carboxylic acid;

3-( {6-[(Methyl-trifluoromethanesulfonyl-amino)-methyl]-naphthalene-2-carbonyl}-amino)-benzoic acid;

1-*tert*-Butoxycarbonylmethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid ethyl ester;

1-Carboxymethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid ethyl ester;

1-*tert*-Butoxycarbonylmethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid;

1-Carboxymethyl-5-(*N,N*-ditrifluoromethanesulfonyl)amino-1*H*-indole-2-carboxylic acid ethyl ester;

1-*tert*-Butoxycarbonylmethyl-5-(*N,N*-ditrifluoromethanesulfonyl)amino-1*H*-indole-2-carboxylic acid ethyl ester;

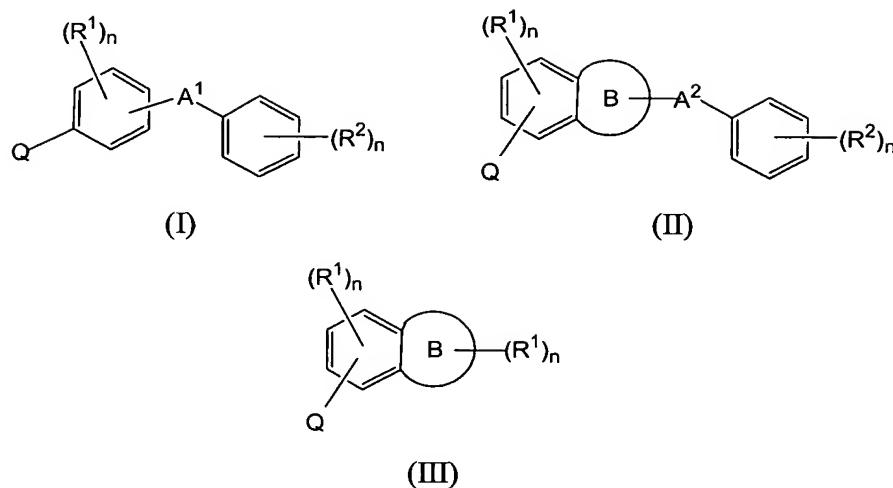
1-Carboxymethyl-5-(*N,N*- ditrifluoromethanesulfonyl)amino-1*H*-indole-2-carboxylic acid;

1-Cyclohexylmethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid ethyl ester;

1-Benzyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid; or

1-Cyclohexylmethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid.

16. (Original) A pharmaceutical composition comprising a compound having the formula:



or a pharmaceutically acceptable salt or solvate thereof, wherein:

$Q$  is  $CF_3SO_2$ ,  $CF_3SO_2NR^3$ ,  $CF_3SO_2R^4$  or  $CF_3SO_2N(R^3)R^4$ , wherein  $R^3$  is H, alkoxy, acyl or  $C_1$ - $C_3$  alkyl, each of which may be substituted or unsubstituted, and  $R^4$  is methylene which may be substituted or unsubstituted;

each  $R^1$  is independently  $C_1$ - $C_3$  alkyl,  $C_1$ - $C_3$  haloalkyl, CN,  $(C=O)OR$ ,  $(C=O)R^5$ , H, halo,  $NHR$ ,  $NH(C=O)OR$ ,  $NH(C=O)R^5$ ,  $NO_2$ ,  $NHSO_2R^5$ ,  $O(C=O)R$ , OR, OH,  $SO_2R^5$ ,  $R^4SO_2CF_3$  or tetrazole, wherein  $R^5$  is  $CF_3$ ,  $C_1$ - $C_3$  alkyl,  $NHR$  and wherein R is H,  $C_1$ - $C_3$  alkyl, aryl or heteroaryl, which may be substituted or unsubstituted;

each  $R^2$  is independently  $C_1$ - $C_3$  alkyl,  $C_1$ - $C_3$  haloalkyl, CN,  $(C=O)OR$ ,  $(C=O)R^5$ , H, halo,  $O(C=O)R$ , OR, OH,  $NHR$ ,  $NH(C=O)OR$ ,  $NH(C=O)R^5$ ,  $NO_2$ ,  $NHSO_2R^5$ ,  $SO_2R^5$ , tetrazole, or  $X_1$ - $R^6$ - $X_2$  wherein  $X_1$  may be present or absent and if present is O, N,  $(C=O)$ ,  $(C=O)NH$ ,  $NH(C=O)$ ,  $SO_2NH$ ,  $NHSO_2$ ;

$R^6$  is  $C_1$ - $3$  alkylene which may be substituted or unsubstituted;

$X_2$  is  $CF_3$ ,  $(C=O)OR$ ,  $(C=O)R^5$ , H,  $NH(C=O)R^5$ ,  $NH(C=O)OR$ ,  $NHSO_2R^5$ ,  $NRR^3$ ,  $O(C=O)R$ , OR,  $SO_2R^5$ , tetrazole;

each  $n$  is independently from 0 to 3;

ring B is an aryl, carbocyclic, heteroaryl, heterocyclic or phenyl ring which may be substituted or unsubstituted;

$A^1$  is a linkage in which the shortest path is 2-8 atoms in length wherein the atoms in the linkage are carbon which may be substituted or unsubstituted or the carbon replaced with

a single nitrogen or oxygen, or combination of nitrogen, oxygen and sulfur provided no two heteroatoms are adjacently linked in a linear linkage; the linkage may be or may contain an aryl, carbocyclic, heteroaryl, heterocyclic or a phenyl ring, which may be directly in the linkage or appended to the linkage; the linkage may be acylalkyl, alkenylene, alkoxy, alkoxyalkyl, alkoxyamino, alkoxyarylalkoxy, alkoxyarylalkyl, alkoxyarylarnino, alkoxyaryloxyalkyl, alkylarnino, alkylarninoalkyl, alkylarninoarylarninoalkyl, alkylaryl, alkylarylalkyl, alkylarylarnino, alkylaryloxy, alkylene, alkylenediamine, alkylenedioxy, alkyloxy, alkyloxyaryl, alkyloxyarylalkyloxy, alkyloxyaryloxyalkyl, C1-C6 alkylsulfonylarnino, alkylthio, alkylthioalkyl, alkynylene, C1-C6 N-sulfonamido, C3-C7 N-amido, aminoalkyl, aminoalkylarnino, aminoalkylarylalkyl, aminoalkylarylalkylarnino, aminoalkylaryloxy, aminoalkyloxy, aminoaryl, aminoarylalkyl, aminoarylcarbonyl, aminoaryloxy, aminoaryloxyalkyl, aminoarylsulfonyl, aryl, arylarnino, aryldioxy, aryldiamine, aryloxy, aryloxyalkyl, aryloxyamino, aryloxyaminoalkyl, aryloxycarbonyl, aryloxysulfonyl, benzimidazole, benzo[b]furan, benzo[b]thiophene, C3-C7 C-amido, carbonylarylarnino, carbonylarylcarbonyl, carbonylaryloxy, chromene, cycloalkylene, disulfide, furan, haloalkyl, imidazole, imidazolidine, imidazoline, indole, isothiazole, isoxazole, morpholine, oxadiazole, oxazole, oxirane, parathiazine, phenothiazine, piperazine, piperidine, purine, pyran, pyrazine, pyrazole, pyrazolidine, pyrimidine, pyridine, pyrrole, pyrrolidine, quinoline, C2-C6 S-sulfonamido, sulfonylalkyl, sulfonylarylarnino, sulfonylaryloxy, sulfonylarylulfonyl, thiadiazole, thiazole, thiophene, triazine, triazole, unsubstituted azeridine, C3-C6 ureido, which may be substituted or unsubstituted;

$A^2$  is a linkage in which the shortest path is 0-6 atoms in length wherein the atoms in the linkage are carbon which may be substituted or unsubstituted or the carbon replaced with a single nitrogen, oxygen or sulfur, or combination of nitrogen, oxygen and sulfur; the linkage may be or may contain an aryl, carbocyclic, heteroaryl, heterocyclic or a phenyl ring, which may be directly in the linkage or appended to the linkage; the linkage may be single atom C, O, S or N which may be substituted or unsubstituted; the linkage may be acylalkyl, alkenylene, alkoxy, alkoxyalkyl, alkoxyamino, alkoxyarylalkoxy, alkoxyarylalkyl, alkoxyarylamino, alkoxyaryloxyalkyl, alkylamino, alkylaminoalkyl, alkylaminoarylaminooalkyl, alkylaryl, alkylarylalkyl, alkylarylamino, alkylaryloxy, alkylene, alkylenediamine, alkylenedioxy, alkyloxy, alkyloxyaryl, alkyloxyarylalkyloxy,

alkyloxyaryloxyalkyl, alkylsulfonylamino, alkylthio, alkylthioalkyl, alkynylene, N-sulfonamido, N-amido, aminoalkyl, aminoalkylamino, aminoalkylarylalkyl, aminoalkylarylalkylamino, aminoalkylaryloxy, aminoalkyloxy, aminoaryl, aminoarylalkyl, aminoarylcarbonyl, aminoaryloxy, aminoaryloxyalkyl, aminoarylsulfonyl, aryl, arylamino, ortho or para aryldioxy, substituted meta-aryldioxy, aryldiamine, aryloxy, aryloxyalkyl, aryloxyamino, aryloxyaminoalkyl, aryloxycarbonyl, aryloxysulfonyl, benzimidazole, benzo[b]furan, benzo[b]thiophene, C-amido, carbonylarylamin, carbonylarylcarbonyl, carbonylaryloxy, chromene, cycloalkylene, furan, haloalkyl, imidazole, imidazolidine, imidazoline, indole, isothiazole, isoxazole, morpholine, oxadiazole, oxazole, oxirane, parathiazine, phenothiazine, piperazine, piperidine, purine, pyran, pyrazine, pyrazole, pyrazolidine, pyrimidine, pyridine, pyrrole, pyrrolidine, quinoline, sulfonamido, sulfonylalkyl, sulfonylarylamin, sulfonylaryloxy, sulfonylarylsulfonyl, thiadiazole, thiazole, thiophene, triazine, triazole, unsubstituted azeridine, ureido, which may be substituted or unsubstituted.

17. (Original) The pharmaceutical composition of claim 16 wherein:

Q is  $\text{CF}_3\text{SO}_2$  or  $\text{CF}_3\text{SO}_2\text{NH}$ ;

each R1 is independently  $\text{CF}_3$ ,  $(\text{C}=\text{O})\text{OR}$ ,  $(\text{C}=\text{O})\text{R}_5$ , H, halo,  $\text{NHR}$ ,  $\text{NH}(\text{C}=\text{O})\text{OR}$ ,  $\text{NH}(\text{C}=\text{O})\text{R}_5$ ,  $\text{NHSO}_2\text{R}_5$ ,  $\text{NO}_2$ ,  $\text{O}(\text{C}=\text{O})\text{R}$ , OH, OR,  $\text{SO}_2\text{R}_5$  or tetrazole;

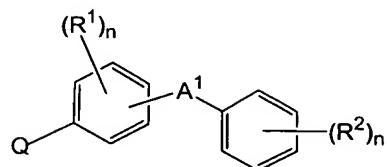
each R2 is independently  $(\text{C}=\text{O})\text{OR}$ ,  $(\text{C}=\text{O})\text{R}_5$ ,  $\text{NH}(\text{C}=\text{O})\text{OR}$ ,  $\text{NH}(\text{C}=\text{O})\text{R}_5$ ,  $\text{NHR}$ ,  $\text{NHSO}_2\text{R}_5$ ,  $\text{NO}_2$ ,  $\text{SO}_2\text{R}_5$ ,  $-\text{R}_6-(\text{C}=\text{O})\text{OR}$ ,  $-\text{R}_6-\text{NRR}_3$ ,  $-\text{R}_6$ -tetrazole or tetrazole;

each n is independently from 0 to 2;

ring B is phenyl or heteroaryl which may be substituted or unsubstituted; and

linkage A1 is C2-C4 alkoxy, C2-C4 alkoxyalkyl, C2-C4 alkylenedioxy, C2-C4 alkylaminoalkyl, C2-C4 alkylenediamine, C3-C4 C-amido, C3-C4 N-amido, C3-C4 ureido, C1-C3 N-sulfonamido, C2-C3 S-sulfonamido, aryldioxy, aryldiamine, aryl, alkylarylalkyl, imidazole, oxazole, oxadiazole, pyrazole, pyrazolidine, pyrrole or triazole.

18. (Original) The pharmaceutical composition of claim 17 wherein the compound has the formula:



(I)

or a pharmaceutically acceptable salt or solvate thereof, wherein:

Q is  $CF_3SO_2$  or  $CF_3SO_2NH$ ;  
 each  $R^1$  is independently H, NHR,  $NO_2$  or OR;  
 each  $R^2$  is independently  $(C=O)OR$ , or  $NHSO_2R^5$  or  $SO_2R^5$ ;  
 each n is independently from 0 to 2; and  
 the linkage  $A^1$  is alkylarylalkyl, C2-C4 alkoxyalkyl, C2-C4 alkylenedioxy, aryl, aryl diamine, aryldioxy or oxadiazole.

19. (Original) The pharmaceutical composition of claim 16 wherein the compound is:

Bis(4-Trifluoromethylsulfonylbenzyl) ether;  
 4-Trifluoromethylsulfonylbenzyl 4-trifluoromethylsulfonylphenyl ether;  
 N,N-Bis(4-trifluoromethylsulfonylbenzyl)benzamide;  
 1,2-Bis(4-trifluoromethylsulfonylphenyl)ethane;  
 N-(4-Trifluoromethylsulfonylbenzyl)-4-trifluoromethylsulfonylbenzamide;  
 N-(4-Trifluoromethylsulfonylbenzyl)benzamide;  
 3,5-Bis-(4-trifluoromethanesulfonyl-phenoxy)-benzoic acid methyl ester;  
 [3,5-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-phenyl]-acetic acid methyl ester;  
 3,5-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-benzoic acid methyl ester;  
 1,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-cyclopentane;  
 4-Methyl-2,6-bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-benzoic acid methyl ester;  
 4-[2-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-ethoxy]-benzoic acid methyl ester;  
 4-[3-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-phenoxy]-benzoic acid;

1-(3,5-Bis-trifluoromethyl-phenyl)-5-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-1*H*-pyrazole-3-carboxylic acid methyl ester;

{4-[4-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-benzenesulfonyl]-phenoxy}-acetic acid ethyl ester;

4-[3-(4-Trifluoromethanesulfonyl-phenoxy)-phenoxy]-benzoic acid;

{4-[4-(4-Trifluoromethanesulfonyl-phenoxy)-benzenesulfonyl]-phenoxy}-acetic acid ethyl ester;

*N*-(3-Trifluoromethanesulfonyl-phenyl)-2-{2-[(3-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;

*N*-(3-Trifluoromethanesulfonyl-phenyl)-2-{3-[(3-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;

*N*-(3-Trifluoromethanesulfonyl-phenyl)-2-{4-[(3-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;

3,6-Bis-(morpholin-4-ylmethyl)-2,5-bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-benzene;

[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-dimethyl-amine;

*N*-(2-Ethylamino-5-trifluoromethanesulfonyl-phenyl)-2-(4-methanesulfonyl-phenyl)-acetamide;

2-Hydroxy-5-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-terephthalic acid diethyl ester;

{2-[(3-Trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetic acid;

{3-[(3-Trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetic acid;

{4-[(3-Trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetic acid;

3,5-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-benzamide;

3,5-Bis-(4-trifluoromethanesulfonyl-phenoxy)-benzoic acid;

*N*-(4-Trifluoromethanesulfonyl-phenyl)-2-{2-[(4-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;

*N*-(4-Trifluoromethanesulfonyl-phenyl)-2-{3-[(4-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;

*N*-(4-Trifluoromethanesulfonyl-phenyl)-2-{4-[(4-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-benzoic acid methyl ester;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-benzoic acid;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-*N*-pyridin-4-yl-benzamide;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-*N*-(4-methoxy-phenyl)-benzamide;

3-[4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-benzoylamino]-benzoic acid ethyl ester;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-*N*-(2-pyrrolidin-1-yl-ethyl)-benzamide;

*N*-Ethyl-4-(1-ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-benzamide;

1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazole-2-carboxylic acid;

[2-(Benzoyl-butyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid methyl ester;

*N*-Benzyl-*N*-[butylcarbamoyl-(4-trifluoromethanesulfonyl-phenyl)-methyl]-benzamide;

*N*-[Butylcarbamoyl-(4-trifluoromethanesulfonyl-phenyl)-methyl]-*N*-(2-hydroxy-ethyl)-benzamide;

[2-(Acetyl-cyclopropyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

[2-(Acetyl-methyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

[2-(Benzoyl-cyclohexyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

*N*-Cyclohexyl-*N*-[(2,6-dimethyl-phenylcarbamoyl)-(4-trifluoromethanesulfonyl-phenyl)-methyl]-benzamide;

[2-(Acetyl-propyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

[2-(Acetyl-cyclohexyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

{4-[4-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-benzenesulfonyl]-phenoxy}-acetic acid;

4-[2-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-ethoxy]-benzoic acid;

2,5-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-terephthalic acid diethyl ester;

1-[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-piperidine;

4-[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-morpholine;

[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-(2-nitro-phenyl)-amine;

1-(2-Nitro-phenylamino)-3-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propan-2-ol;

[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-(4-nitro-phenyl)-amine;

1-(4-Nitro-phenylamino)-3-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propan-2-ol;

4-[2-Hydroxy-3-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propylamino]-benzenesulfonamide; or

4-[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propylamino]-benzenesulfonamide.

20. (Original) The pharmaceutical composition of claim 16 wherein the compound is:

1,2-Bis(4-trifluoromethylsulfonamidophenyl)ethane;

1,2-Bis(2-methyl-4-trifluoromethylsulfonamidophenyl)ethane;

1,3-Bis(4-trifluoromethylsulfonamidophenoxy)-2,2-dimethylpropane;

1,3-Bis(4-trifluoromethylsulfonamidophenoxy)propane;

1,4-Bis(4-trifluoromethylsulfonamidophenoxy)butane;

1,4-Bis(4-trifluoromethylsulfonamidophenoxy)benzene;

1-(4-Aminophenoxy)-4-trifluoromethylsulfonamidophenoxy benzene;

Bis(4-trifluoromethylsulfonamidophenyl) ether;

1,3-Bis(4-trifluoromethylsulfonamidophenoxy)benzene;

2,5-Bis(4-trifluoromethylsulfonamidophenyl)-(1,3,4)oxadiazole;

Bis(4-trifluoromethylsulfonamidophenyl)-1,4-diisopropylbenzene;  
5-Trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid;  
1-Methyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid;  
(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetic acid;  
1-Methyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid phenylamide;  
5-Trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid phenylamide;  
3-[(1-Methyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carbonyl)-amino]-benzoic acid;  
3-[(5-Trifluoromethanesulfonylamino-1*H*-indole-2-carbonyl)-amino]-benzoic acid;  
4-[(5-Trifluoromethanesulfonylamino-1*H*-indole-2-carbonyl)-amino]-benzoic acid;  
4-[2-(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetylamino]-benzoic acid;  
3-[2-(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetylamino]-benzoic acid;  
4-{{[2-(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetylamino]-methyl}-benzoic acid;  
(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetic acid *tert*-butyl ester;  
1-Methyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid ethyl ester;  
6-Trifluoromethanesulfonylamino-naphthalene-2-carboxylic acid;  
*N,N*-Bis[(6-carboxyl-naphthalen-2-yl)methyl] trifluoromethanesulfonamide;  
6-[(Methyl-trifluoromethanesulfonyl-amino)-methyl]-naphthalene-2-carboxylic acid;  
3-({6-[(Methyl-trifluoromethanesulfonyl-amino)-methyl]-naphthalene-2-carbonyl}-amino)-benzoic acid;  
1-*tert*-Butoxycarbonylmethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid ethyl ester;  
1-Carboxymethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid ethyl ester;  
1-*tert*-Butoxycarbonylmethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid;  
1-Carboxymethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid;

1-Carboxymethyl-5-(*N,N*-ditrifluoromethanesulfonyl)amino-1*H*-indole-2-carboxylic acid ethyl ester;

1-*tert*-Butoxycarbonylmethyl-5-(*N,N*-ditrifluoromethanesulfonyl)amino-1*H*-indole-2-carboxylic acid ethyl ester;

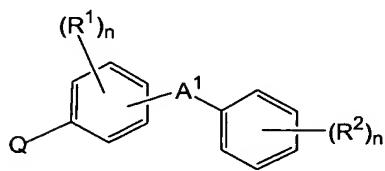
1-Carboxymethyl-5-(*N,N*-ditrifluoromethanesulfonyl)amino-1*H*-indole-2-carboxylic acid;

1-Cyclohexylmethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid ethyl ester;

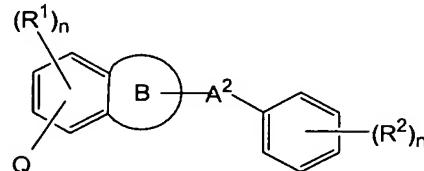
1-Benzyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid; or

1-Cyclohexylmethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid.

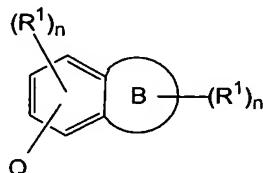
21. (Original) A method for treating a protein tyrosine phosphatase signal transduction associated disorder in a mammal which comprises administering to the mammal therapeutically effective amount of a compound having the formula:



(I)



(II)



(III)

or a pharmaceutically acceptable salt or solvate thereof, wherein:

Q is  $\text{CF}_3\text{SO}_2$ ,  $\text{CF}_3\text{SO}_2\text{NR}^3$ ,  $\text{CF}_3\text{SO}_2\text{R}^4$  or  $\text{CF}_3\text{SO}_2\text{N}(\text{R}^3)\text{R}^4$ , wherein  $\text{R}^3$  is H, alkoxy, acyl or  $\text{C}_1\text{-C}_3$  alkyl, each of which may be substituted or unsubstituted, and  $\text{R}^4$  is methylene which may be substituted or unsubstituted;

each  $\text{R}^1$  is independently  $\text{C}_1\text{-C}_3$  alkyl,  $\text{C}_1\text{-C}_3$  haloalkyl, CN,  $(\text{C}=\text{O})\text{OR}$ ,  $(\text{C}=\text{O})\text{R}^5$ , H, halo, NHR,  $\text{NH}(\text{C}=\text{O})\text{OR}$ ,  $\text{NH}(\text{C}=\text{O})\text{R}^5$ , NO<sub>2</sub>,  $\text{NHSO}_2\text{R}^5$ ,  $\text{O}(\text{C}=\text{O})\text{R}$ , OH, OR, SO<sub>2</sub>R<sub>5</sub>,

R4SO2CF3 or tetrazole, wherein R5 is CF3, C1-C3 alkyl, NHR and wherein R is H, C1-C3 alkyl, aryl or heteroaryl, which may be substituted or unsubstituted;

each R2 is independently C1-C3 alkyl, C1-C3 haloalkyl, CN, (C=O)OR, (C=O)R5, H, halo, O(C=O)R, OR, OH, NHR, NH(C=O)OR, NH(C=O)R5, NO2, NSO2R5, SO2R5, tetrazole, or X1-R6-X2, wherein X1 may be present or absent and if present is O, N, (C=O), (C=O)NH, NH(C=O), SO2NH, NSO2, R6 is C1-3 alkylene which may be substituted or unsubstituted and X2 is CF3, (C=O)OR, (C=O)R5, H, NH(C=O)R5, NH(C=O)OR, NSO2R5, NRR3, O(C=O)R, OR, SO2R5, tetrazole;

each n is independently from 0 to 3;

ring B is an aryl, carbocyclic, heteroaryl, heterocyclic or phenyl ring which may be substituted or unsubstituted;

A1 is a linkage in which the shortest path is 2-8 atoms in length wherein the atoms in the linkage are carbon which may be substituted or unsubstituted or the carbon replaced with a single nitrogen, oxygen or sulfur, or combination of nitrogen, oxygen and sulfur; the linkage may be or may contain an aryl, carbocyclic, heteroaryl, heterocyclic or a phenyl ring, which may be directly in the linkage or appended to the linkage; the linkage may be acylalkyl, alkenylene, alkoxy, alkoxyalkyl, alkoxyamino, alkoxyarylalkoxy, alkoxyarylalkyl, alkoxyarylamino, alkoxyaryloxyalkyl, alkylamino, alkylaminoalkyl, alkylaminoarylaminolalkyl, alkylaryl, alkylarylalkyl, alkylarylamino, alkylaryloxy, alkylene, alkylenediamine, alkylenedioxy, alkyloxy, alkyloxyaryl, alkyloxyarylalkyloxy, alkyloxyaryloxyalkyl, alkylsulfonylamino, alkylthio, alkylthioalkyl, alkynylene, N-sulfonamido, N-amido, aminoalkyl, aminoalkylamino, aminoalkylarylalkyl, aminoalkylarylalkylamino, aminoalkylaryloxy, aminoalkyloxy, aminoaryl, aminoarylalkyl, aminoarylcarbonyl, aminoaryloxy, aminoaryloxyalkyl, aminoarylsulfonyl, aryl, arylamino, aryldioxy, aryldiamine, aryloxy, aryloxyalkyl, aryloxyamino, aryloxyaminoalkyl, aryloxycarbonyl, aryloxysulfonyl, benzimidazole, benzo[b]furan, benzo[b]thiophene, C-amido, carbonylarylaminol, carbonylarylcarbonyl, carbonylaryloxy, chromene, cycloalkylene, disulfide, furan, haloalkyl, imidazole, imidazolidine, imidazoline, indole, isothiazole, isoxazole, morpholine, oxadiazole, oxazole, oxirane, parathiazine, phenothiazine, piperazine, piperidine, purine, pyran, pyrazine, pyrazole, pyrazolidine, pyrimidine, pyridine, pyrrole, pyrrolidine, quinoline, sulfonamido, sulfonylalkyl, sulfonylarylaminol, sulfonylaryloxy,

sulfonylarylsulfonyl, thiadiazole, thiazole, thiophene, triazine, triazole, unsubstituted azeridine, ureido, which may be substituted or unsubstituted;

$A^2$  is a linkage in which the shortest path is 0-6 atoms in length wherein the atoms in the linkage are carbon which may be substituted or unsubstituted or the carbon replaced with a single nitrogen, oxygen or sulfur, or combination of nitrogen, oxygen and sulfur; the linkage may be or may contain an aryl, carbocyclic, heteroaryl, heterocyclic or a phenyl ring, which may be directly in the linkage or appended to the linkage; the linkage may be single atom C, O, S or N which may be substituted or unsubstituted; the linkage may be acylalkyl, alkenylene, alkoxy, alkoxyalkyl, alkoxyamino, alkoxyarylalkoxy, alkoxyarylalkyl, alkoxyarylamino, alkoxyaryloxyalkyl, alkylamino, alkylaminoalkyl, alkylaminoarylaminooalkyl, alkylaryl, alkylarylalkyl, alkylarylamino, alkylaryloxy, alkylene, alkylenediamine, alkylenedioxy, alkyloxy, alkyloxyaryl, alkyloxyarylalkyloxy, alkyloxyaryloxyalkyl, alkylsulfonylamino, alkylthio, alkylthioalkyl, alkynylene, N-sulfonamido, N-amido, aminoalkyl, aminoalkylamino, aminoalkylarylalkyl, aminoalkylarylalkylamino, aminoalkylaryloxy, aminoalkyloxy, aminoaryl, aminoarylalkyl, aminoarylcarbonyl, aminoaryloxy, aminoaryloxyalkyl, aminoarylsulfonyl, aryl, arylamino, ortho or para aryldioxy, substituted meta-aryldioxy, arylamine, aryloxy, aryloxyalkyl, aryloxyamino, aryloxyaminoalkyl, aryloxycarbonyl, aryloxysulfonyl, benzimidazole, benzo[b]furan, benzo[b]thiophene, C-amido, carbonylarylaminoo, carbonylarylcarbonyl, carbonylaryloxy, chromene, cycloalkylene, disulfide, furan, haloalkyl, imidazole, imidazolidine, imidazoline, indole, isothiazole, isoxazole, morpholine, oxadiazole, oxazole, oxirane, parathiazine, phenothiazine, piperazine, piperidine, purine, pyran, pyrazine, pyrazole, pyrazolidine, pyrimidine, pyridine, pyrrole, pyrrolidine, quinoline, sulfonamido, sulfonylalkyl, sulfonylarylaminoo, sulfonylaryloxy, sulfonylarylsulfonyl, thiadiazole, thiazole, thiophene, triazine, triazole, unsubstituted azeridine, ureido, which may be substituted or unsubstituted

and a pharmaceutically acceptable carrier or excipient;

wherein said compound regulates, inhibits or modulates protein tyrosine phosphatase signal transduction.

22. (Original) The method of claim 21 wherein the protein tyrosine phosphatase signal transduction is associated with cancer, a solid tumor, glioma, melanoma, Kaposi's sarcoma, hemangioma, ovarian cancer, breast cancer, lung cancer, pancreatic cancer, liver cancer, prostate cancer, colon cancer, or epidermoid cancer.

23. (Original) The method of claim 21 wherein the protein tyrosine phosphatase signal transduction is associated with diabetes.

24. (Original) The method of claim 21 wherein the protein tyrosine phosphatase signal transduction is associated with neurological degenerative diseases.

25. (Original) The method of claim 21 wherein the protein tyrosine phosphatase signal transduction is associated with osteoporosis.

26. (Currently Amended) The method of claim 21, 21, 22, 23, 24 or 25 wherein the mammal is a human.

27. (Original) The method of claim 21 wherein Q is  $\text{CF}_3\text{SO}_2$ .

28. (Original) The method of claim 21 wherein Q is  $\text{CF}_3\text{SO}_2\text{NR}^3$ .

29. (Original) The method of claim 21 wherein Q is  $\text{CF}_3\text{SO}_2\text{R}^4$ .

30. (Original) The method of claim 21 wherein Q is  $\text{CF}_3\text{SO}_2\text{N}(\text{R}^3)\text{R}^4$ .

31. (Original) The method of claim 21 wherein the compound has formula I.

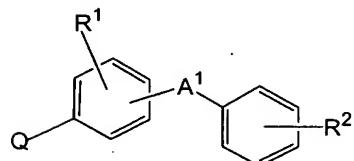
32. (Original) The method of claim 21 wherein the compound has formula II.

33. (Original) The method of claim 21 wherein the compound has formula III.

34. (Original) The method of claim 21 wherein R<sup>2</sup> is SO<sub>2</sub>R<sup>5</sup>, NHSO<sub>2</sub>R<sup>5</sup> or CF<sub>3</sub>SO<sub>2</sub>R<sup>4</sup>.

35. (Original) The method of claim 21 wherein Q is CF<sub>3</sub>SO<sub>2</sub> or CF<sub>3</sub>SO<sub>2</sub>NH; each R1 is independently CF<sub>3</sub>, (C=O)OR, (C=O)R5, H, halo, NHR, NH(C=O)OR, NH(C=O)R5, NHSO<sub>2</sub>R5, NO<sub>2</sub>, O(C=O)R, OH, OR, SO<sub>2</sub>R5 or tetrazole; each R2 is independently (C=O)OR, (C=O)R5, NH(C=O)OR, NH(C=O)R5, NHR, NHSO<sub>2</sub>R5, NO<sub>2</sub>, SO<sub>2</sub>R5, -R6-(C=O)OR, -R6-NRR3, -R6-tetrazole or tetrazole; each n is independently from 0 to 2; ring B is phenyl or heteroaryl which may be substituted or unsubstituted; and linkage A1 is C1-C4 alkoxy, C2-C4 alkoxyalkyl, C2-C4 alkylenedioxy, C2-C4 alkylaminoalkyl, C2-C4 alkylenediamine, C1-C4 C-amido, C1-C4 N-amido, C1-C4 ureido, C0-C3 N-sulfonamido, C0-C3 S-sulfonamido, aryldioxy, aryldiamine, aryl, alkylarylalkyl, imidazole, oxazole, oxadiazole, pyrazole, pyrazolidine, pyrrole or triazole.

36. (Original) The method of claim 21 wherein the compound has the formula:



(IV)

or a pharmaceutically acceptable salt or solvate thereof, wherein:

Q is CF<sub>3</sub>SO<sub>2</sub> or CF<sub>3</sub>SO<sub>2</sub>NH;

R1 is H or NO<sub>2</sub>;

R2 is (C=O)OR, NHSO<sub>2</sub>R5 or SO<sub>2</sub>R5; and

the linkage A1 is C2-C4 alkoxyalkyl, aryldioxy, aryl, alkylarylalkyl, C1-C4 N-amido or oxadiazole.

37. (Original) The method of claim 21 wherein the compound is:

Bis(4-Trifluoromethylsulfonylbenzyl) ether;  
4-Trifluoromethylsulfonylbenzyl 4-trifluoromethylsulfonylphenyl ether;  
N,N-Bis(4-trifluoromethylsulfonylbenzyl)benzamide;  
1,2-Bis(4-trifluoromethylsulfonylphenyl)ethane;  
N-(4-Trifluoromethylsulfonylbenzyl)-4-trifluoromethylsulfonylbenzamide;  
N-(4-Trifluoromethylsulfonylbenzyl)benzamide;  
Bis(4-Trifluoromethylsulfonylphenyl) disulfide;  
Bis-(2-Nitro-4-trifluoromethylsulfonylphenyl) disulfide;  
3,5-Bis-(4-trifluoromethanesulfonyl-phenoxy)-benzoic acid methyl ester;  
[3,5-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-phenyl]-acetic acid methyl ester;  
3,5-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-benzoic acid methyl ester;  
1,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-cyclopentane;  
4-Methyl-2,6-bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-benzoic acid methyl ester;  
4-[2-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-ethoxy]-benzoic acid methyl ester;  
4-[3-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-phenoxy]-benzoic acid;  
1-(3,5-Bis-trifluoromethyl-phenyl)-5-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-1*H*-pyrazole-3-carboxylic acid methyl ester;  
{4-[4-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-benzenesulfonyl]-phenoxy}-acetic acid ethyl ester;  
4-[3-(4-Trifluoromethanesulfonyl-phenoxy)-phenoxy]-benzoic acid;  
{4-[4-(4-Trifluoromethanesulfonyl-phenoxy)-benzenesulfonyl]-phenoxy}-acetic acid ethyl ester;  
N-(3-Trifluoromethanesulfonyl-phenyl)-2-{2-[(3-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;  
N-(3-Trifluoromethanesulfonyl-phenyl)-2-{3-[(3-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;  
N-(3-Trifluoromethanesulfonyl-phenyl)-2-{4-[(3-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;

3,6-Bis-(morpholin-4-ylmethyl)-2,5-bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-benzene;

[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-dimethyl-amine;

*N*-(2-Ethylamino-5-trifluoromethanesulfonyl-phenyl)-2-(4-methanesulfonyl-phenyl)-acetamide;

2-Hydroxy-5-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-terephthalic acid diethyl ester;

{2-[(3-Trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetic acid;

{3-[(3-Trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetic acid;

{4-[(3-Trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetic acid;

3,5-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-benzamide;

3,5-Bis-(4-trifluoromethanesulfonyl-phenoxy)-benzoic acid;

*N*-(4-Trifluoromethanesulfonyl-phenyl)-2-{2-[(4-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;

*N*-(4-Trifluoromethanesulfonyl-phenyl)-2-{3-[(4-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;

*N*-(4-Trifluoromethanesulfonyl-phenyl)-2-{4-[(4-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-benzoic acid methyl ester;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-benzoic acid;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-*N*-pyridin-4-yl-benzamide;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-*N*-(4-methoxy-phenyl)-benzamide;

3-[4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-benzoylamino]-benzoic acid ethyl ester;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-*N*-(2-pyrrolidin-1-yl-ethyl)-benzamide;

*N*-Ethyl-4-(1-ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-benzamide;

1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazole-2-carboxylic acid;

[2-(Benzoyl-butyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid methyl ester;

*N*-Benzyl-*N*-[butylcarbamoyl-(4-trifluoromethanesulfonyl-phenyl)-methyl]-benzamide;

*N*-[Butylcarbamoyl-(4-trifluoromethanesulfonyl-phenyl)-methyl]-*N*-(2-hydroxyethyl)-benzamide;

[2-(Acetyl-cyclopropyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

[2-(Acetyl-methyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

[2-(Benzoyl-cyclohexyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

*N*-Cyclohexyl-*N*-[(2,6-dimethyl-phenylcarbamoyl)-(4-trifluoromethanesulfonyl-phenyl)-methyl]-benzamide;

[2-(Acetyl-propyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

[2-(Acetyl-cyclohexyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

{4-[4-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-benzenesulfonyl]-phenoxy}-acetic acid;

4-[2-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-ethoxy]-benzoic acid;

2,5-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-terephthalic acid diethyl ester;

1-[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-piperidine;

4-[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-morpholine;

[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-(2-nitro-phenyl)-amine;

1-(2-Nitro-phenylamino)-3-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propan-2-ol;

[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-(4-nitro-phenyl)-amine;

1-(4-Nitro-phenylamino)-3-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propan-2-ol;

4-[2-Hydroxy-3-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propylamino]-benzenesulfonamide;

4-[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propylamino]-benzenesulfonamide; or

Bis- {[4-(2-nitro-4-trifluoromethanesulfonyl)-phenoxy]-phenyl} sulfone.

38. (Original) The method of claim 21 wherein the compound is:

1,2-Bis(4-trifluoromethylsulfonamidophenyl)ethane;

1,2-Bis(2-methyl-4-trifluoromethylsulfonamidophenyl)ethane;

1,3-Bis(4-trifluoromethylsulfonamidophenoxy)-2,2-dimethylpropane;

1,3-Bis(4-trifluoromethylsulfonamidophenoxy)propane;

1,4-Bis(4-trifluoromethylsulfonamidophenoxy)butane;

1,4-Bis(4-trifluoromethylsulfonamidophenoxy)benzene;

1-(4-Aminophenoxy)-4-trifluoromethylsulfonamidophenoxy benzene;

Bis(4-trifluoromethylsulfonamidophenyl) ether;

1,3-Bis(4-trifluoromethylsulfonamidophenoxy)benzene;

2,5-Bis(4-trifluoromethylsulfonamidophenyl)-(1,3,4)oxadiazole;

Bis(4-trifluoromethylsulfonamidophenyl)-1,4-diisopropylbenzene;

5-Trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid;

1-Methyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid;

(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetic acid;

1-Methyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid phenylamide;

5-Trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid phenylamide;

3-[(1-Methyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carbonyl)-amino]-benzoic acid;

3-[(5-Trifluoromethanesulfonylamino-1*H*-indole-2-carbonyl)-amino]-benzoic acid;

4-[(5-Trifluoromethanesulfonylamino-1*H*-indole-2-carbonyl)-amino]-benzoic acid;

4-[2-(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetylamino]-benzoic acid;

3-[2-(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetylamino]-benzoic acid;

4- {[2-(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetylamino]-methyl}-benzoic acid;

(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetic acid *tert*-butyl ester;

1-Methyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid ethyl ester;

6-Trifluoromethanesulfonylamino-naphthalene-2-carboxylic acid;

*N,N*-Bis[(6-carboxyl-naphthalen-2-yl)methyl] trifluoromethanesulfonamide;

6-[(Methyl-trifluoromethanesulfonyl-amino)-methyl]-naphthalene-2-carboxylic acid;

3-( {6-[(Methyl-trifluoromethanesulfonyl-amino)-methyl]-naphthalene-2-carbonyl}-amino)-benzoic acid;

1-*tert*-Butoxycarbonylmethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid ethyl ester;

1-Carboxymethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid ethyl ester;

1-*tert*-Butoxycarbonylmethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid;

1-Carboxymethyl-5-(*N,N*-ditrifluoromethanesulfonyl)amino-1*H*-indole-2-carboxylic acid ethyl ester;

1-*tert*-Butoxycarbonylmethyl-5-(*N,N*-ditrifluoromethanesulfonyl)amino-1*H*-indole-2-carboxylic acid ethyl ester;

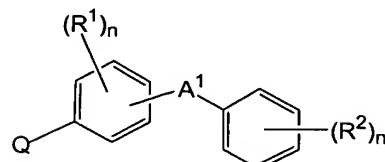
1-Carboxymethyl-5-(*N,N*- ditrifluoromethanesulfonyl)amino-1*H*-indole-2-carboxylic acid;

1-Cyclohexylmethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid ethyl ester;

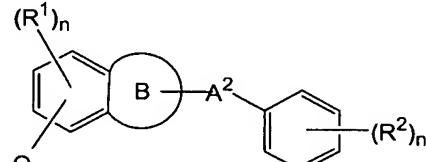
1-Benzyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid; or

1-Cyclohexylmethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid.

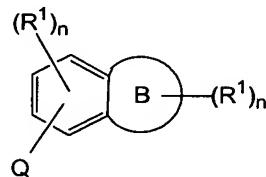
39. (Original) A method for treating, alleviating or preventing cancer in a mammal which comprises administering to a mammal in need thereof a therapeutically effective amount of a compound having the formula:



(I)



(II)



(III)

or a pharmaceutically acceptable salt or solvate thereof, wherein:

Q is  $\text{CF}_3\text{SO}_2$ ,  $\text{CF}_3\text{SO}_2\text{NR}^3$ ,  $\text{CF}_3\text{SO}_2\text{R}^4$  or  $\text{CF}_3\text{SO}_2\text{N}(\text{R}^3)\text{R}^4$ , wherein  $\text{R}^3$  is H, alkoxy, acyl or  $\text{C}_1\text{-C}_3$  alkyl, each of which may be substituted or unsubstituted, and  $\text{R}^4$  is methylene which may be substituted or unsubstituted;

each  $\text{R}1$  is independently  $\text{C}_1\text{-C}_3$  alkyl,  $\text{C}_1\text{-C}_3$  haloalkyl, CN,  $(\text{C}=\text{O})\text{OR}$ ,  $(\text{C}=\text{O})\text{R}5$ , H, halo,  $\text{NHR}$ ,  $\text{NH}(\text{C}=\text{O})\text{OR}$ ,  $\text{NH}(\text{C}=\text{O})\text{R}5$ ,  $\text{NO}_2$ ,  $\text{NHSO}_2\text{R}5$ ,  $\text{O}(\text{C}=\text{O})\text{R}$ , OR, OH,  $\text{SO}_2\text{R}5$ ,  $\text{R}4\text{SO}_2\text{CF}_3$  or tetrazole, wherein  $\text{R}5$  is  $\text{CF}_3$ ,  $\text{C}_1\text{-C}_3$  alkyl,  $\text{NHR}$  and wherein R is H,  $\text{C}_1\text{-C}_3$  alkyl, aryl or heteroaryl, which may be substituted or unsubstituted;

each  $\text{R}2$  is independently  $\text{C}_1\text{-3}$  alkyl,  $\text{C}_1\text{-C}_3$  haloalkyl, CN,  $(\text{C}=\text{O})\text{OR}$ ,  $(\text{C}=\text{O})\text{R}5$ , H, halo,  $\text{O}(\text{C}=\text{O})\text{R}$ , OR, OH,  $\text{NHR}$ ,  $\text{NH}(\text{C}=\text{O})\text{OR}$ ,  $\text{NH}(\text{C}=\text{O})\text{R}5$ ,  $\text{NO}_2$ ,  $\text{NHSO}_2\text{R}5$ ,  $\text{SO}_2\text{R}5$ , tetrazole or  $\text{X}_1\text{-R}_6\text{-X}_2$ , wherein  $\text{X}_1$  may be present or absent and if present is O, N,  $(\text{C}=\text{O})$ ,  $(\text{C}=\text{O})\text{NH}$ ,  $\text{NH}(\text{C}=\text{O})$ ,  $\text{SO}_2\text{NH}$ ,  $\text{NHSO}_2$ ,  $\text{R}_6$  is  $\text{C}_1\text{-3}$  alkylene which may be substituted or unsubstituted and  $\text{X}_2$  is  $\text{CF}_3$ ,  $(\text{C}=\text{O})\text{OR}$ ,  $(\text{C}=\text{O})\text{R}5$ , H,  $\text{NH}(\text{C}=\text{O})\text{R}5$ ,  $\text{NH}(\text{C}=\text{O})\text{OR}$ ,  $\text{NHSO}_2\text{R}5$ ,  $\text{NRR}3$ ,  $\text{O}(\text{C}=\text{O})\text{R}$ , OR,  $\text{SO}_2\text{R}5$ , tetrazole;

each  $n$  is independently from 0 to 3;

ring B is an aryl, carbocyclic, heteroaryl, heterocyclic or phenyl ring which may be substituted or unsubstituted;

A1 is a linkage in which the shortest path is 2-8 atoms in length wherein the atoms in the linkage are carbon which may be substituted or unsubstituted or the carbon replaced with a single nitrogen, oxygen, sulfur, or combination of nitrogen, oxygen and sulfur; the linkage may be or contain an aryl, carbocyclic, heteroaryl, heterocyclic or a phenyl ring, which may be directly in the linkage or appended to the linkage; the linkage may be acylalkyl, alkenylene, alkoxy, alkoxyalkyl, alkoxyamino, alkoxyarylalkoxy, alkoxyarylalkyl, alkoxyarylamino, alkoxyaryloxyalkyl, alkylamino, alkylaminoalkyl, alkylaminoarylaminolalkyl, alkylaryl, alkylarylalkyl, alkylarylamino, alkylaryloxy, alkylene, alkylenediamine, alkylenedioxy, alkyloxy, alkyloxyaryl, alkyloxyarylalkoxy, alkyloxyaryloxyalkyl, alkylsulfonylaminol, alkylthio, alkylthioalkyl, alkynylene, N-sulfonamido, N-amido, aminoalkyl, aminoalkylamino, aminoalkylarylalkyl, aminoalkylarylalkylamino, aminoalkylaryloxy, aminoalkyloxy, aminoaryl, aminoarylalkyl, aminoarylcarbonyl, aminoaryloxy, aminoaryloxyalkyl, aminoarylsulfonyl, aryl, arylamino, aryldioxy, aryldiamine, aryloxy, aryloxyalkyl, aryloxyamino, aryloxyaminoalkyl, aryloxycarbonyl, aryloxysulfonyl, benzimidazole, benzo[b]furan, benzo[b]thiophene, C-amido, carbonylarylaminol, carbonylarylcarbonyl, carbonylaryloxy, chromene, cycloalkylene, disulfide, furan, haloalkyl, imidazole, imidazolidine, imidazoline, indole, isothiazole, isoxazole, morpholine, oxadiazole, oxazole, oxirane, parathiazine, phenothiazine, piperazine, piperidine, purine, pyran, pyrazine, pyrazole, pyrazolidine, pyrimidine, pyridine, pyrrole, pyrrolidine, quinoline, sulfonamido, sulfonylalkyl, sulfonylarylaminol, sulfonylaryloxy, sulfonylarylsulfonyl, thiadiazole, thiazole, thiophene, triazine, triazole, unsubstituted azeridine, or ureido which may be substituted or unsubstituted;

A<sup>2</sup> is a linkage in which the shortest path is 0-6 atoms in length wherein the atoms in the linkage are carbon which may be substituted or unsubstituted or the carbon replaced with a single nitrogen, oxygen or sulfur, or combination of nitrogen, oxygen and sulfur; the linkage may be or may contain an aryl, carbocyclic, heteroaryl, heterocyclic or a phenyl ring, which may be directly in the linkage or appended to the linkage; the linkage may be single atom C, O, S or N which may be substituted or unsubstituted; the linkage may be acylalkyl, alkenylene, alkoxy, alkoxyalkyl, alkoxyamino, alkoxyarylalkoxy, alkoxyarylalkyl, alkoxyarylamino, alkoxyaryloxyalkyl, alkylamino, alkylaminoalkyl, alkylaminoarylaminolalkyl, alkylaryl, alkylarylalkyl, alkylarylamino, alkylaryloxy, alkylene,

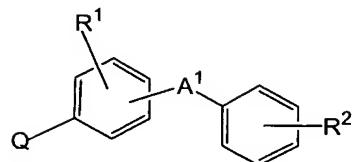
alkylenediamine, alkyleneoxy, alkyloxy, alkyloxyaryl, alkyloxyarylalkyloxy, alkyloxyaryloxyalkyl, alkylsulfonylamino, alkylthio, alkylthioalkyl, alkynylene, N-sulfonamido, N-amido, aminoalkyl, aminoalkylamino, aminoalkylarylalkyl, aminoalkylarylalkylamino, aminoalkylaryloxy, aminoalkyloxy, aminoaryl, aminoarylalkyl, aminoarylcarbonyl, aminoaryloxy, aminoaryloxyalkyl, aminoarylsulfonyl, aryl, arylamino, ortho or para aryldioxy, substituted meta-aryldioxy, aryldiamine, aryloxy, aryloxyalkyl, aryloxyamino, aryloxyaminoalkyl, aryloxycarbonyl, aryloxysulfonyl, benzimidazole, benzo[b]furan, benzo[b]thiophene, C-amido, carbonylarylalmino, carbonylarylcarbonyl, carbonylaryloxy, chromene, cycloalkylene, furan, haloalkyl, imidazole, imidazolidine, imidazoline, indole, isothiazole, isoxazole, morpholine, oxadiazole, oxazole, oxirane, parathiazine, phenothiazine, piperazine, piperidine, purine, pyran, pyrazine, pyrazole, pyrazolidine, pyrimidine, pyridine, pyrrole, pyrrolidine, quinoline, sulfonamido, sulfonylalkyl, sulfonylarylalmino, sulfonylaryloxy, sulfonylarylsulfonyl, thiadiazole, thiazole, thiophene, triazine, triazole, unsubstituted azeridine, ureido, which may be substituted or unsubstituted;

and a pharmaceutically acceptable carrier or excipient.

40. (Original) The method of claim 39 wherein the mammal is a human.

41. (Original) The method of claim 39 wherein  
Q is  $\text{CF}_3\text{SO}_2$  or  $\text{CF}_3\text{SO}_2\text{NH}$ ;  
each R1 is independently  $\text{CF}_3$ ,  $(\text{C}=\text{O})\text{OR}$ ,  $(\text{C}=\text{O})\text{R}_5$ , H, halo, NHR,  $\text{NH}(\text{C}=\text{O})\text{OR}$ ,  $\text{NH}(\text{C}=\text{O})\text{R}_5$ ,  $\text{NHSO}_2\text{R}_5$ ,  $\text{NO}_2$ ,  $\text{O}(\text{C}=\text{O})\text{R}$ , OH, OR,  $\text{SO}_2\text{R}_5$  or tetrazole;  
each R2 is independently  $(\text{C}=\text{O})\text{OR}$ ,  $(\text{C}=\text{O})\text{R}_5$ ,  $\text{NH}(\text{C}=\text{O})\text{OR}$ ,  $\text{NH}(\text{C}=\text{O})\text{R}_5$ , NHR,  $\text{NHSO}_2\text{R}_5$ ,  $\text{NO}_2$ ,  $\text{SO}_2\text{R}_5$ ,  $-\text{R}_6-(\text{C}=\text{O})\text{OR}$ ,  $-\text{R}_6-\text{NRR}_3$ ,  $-\text{R}_6$ -tetrazole or tetrazole;  
each n is independently from 0 to 2;  
ring B is phenyl or heteroaryl which may be substituted or unsubstituted; and  
linkage A1 is C1-C4 alkoxy, C2-C4 alkoxyalkyl, C2-C4 alkyleneoxy, C2-C4 alkylaminoalkyl, C2-C4 alkylene diamine, C1-C4 C-amido, C1-C4 N-amido, C1-C4 ureido, C0-C3 N-sulfonamido, C0-C3 S-sulfonamido, aryldioxy, aryldiamine, aryl, alkylarylalkyl, imidazole, oxazole, oxadiazole, pyrazole, pyrazolidine, pyrrole or triazole.

42. (Original) The method of claim 39 wherein the compound has the formula:



(IV)

or a pharmaceutically acceptable salt or solvate thereof, wherein:

$Q$  is  $CF_3SO_2$  or  $CF_3SO_2NH$ ;

$R^1$  is H or  $NO_2$ ;

$R^2$  is  $(C=O)OR$ ,  $NHSO_2R^5$  or  $SO_2R^5$ ; and

the linkage  $A^1$  is C2-C4 alkoxyalkyl, aryldioxy, aryl, alkylarylkyl, C1-C4 N-amido or oxadiazole.

43. (Original) The method of claim 39 wherein the compound is:

Bis(4-Trifluoromethylsulfonylbenzyl) ether;

4-Trifluoromethylsulfonylbenzyl 4-trifluoromethylsulfonylphenyl ether;

N,N-Bis(4-trifluoromethylsulfonylbenzyl)benzamide;

1,2-Bis(4-trifluoromethylsulfonylphenyl)ethane;

N-(4-Trifluoromethylsulfonylbenzyl)-4-trifluoromethylsulfonylbenzamide;

N-(4-Trifluoromethylsulfonylbenzyl)benzamide;

Bis(4-Trifluoromethylsulfonylphenyl) disulfide;

Bis-(2-Nitro-4-trifluoromethylsulfonylphenyl) disulfide;

3,5-Bis-(4-trifluoromethanesulfonyl-phenoxy)-benzoic acid methyl ester;

[3,5-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-phenyl]-acetic acid methyl ester;

3,5-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-benzoic acid methyl ester;

1,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-cyclopentane;

4-Methyl-2,6-bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-benzoic acid methyl ester;

4-[2-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-ethoxy]-benzoic acid methyl ester;  
4-[3-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-phenoxy]-benzoic acid;  
1-(3,5-Bis-trifluoromethyl-phenyl)-5-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-  
1*H*-pyrazole-3-carboxylic acid methyl ester;  
{4-[4-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-benzenesulfonyl]-phenoxy}-  
acetic acid ethyl ester;  
4-[3-(4-Trifluoromethanesulfonyl-phenoxy)-phenoxy]-benzoic acid;  
{4-[4-(4-Trifluoromethanesulfonyl-phenoxy)-benzenesulfonyl]-phenoxy}-acetic acid  
ethyl ester;  
*N*-(3-Trifluoromethanesulfonyl-phenyl)-2-{2-[(3-trifluoromethanesulfonyl-  
phenylcarbamoyl)-methyl]-phenyl}-acetamide;  
*N*-(3-Trifluoromethanesulfonyl-phenyl)-2-{3-[(3-trifluoromethanesulfonyl-  
phenylcarbamoyl)-methyl]-phenyl}-acetamide;  
*N*-(3-Trifluoromethanesulfonyl-phenyl)-2-{4-[(3-trifluoromethanesulfonyl-  
phenylcarbamoyl)-methyl]-phenyl}-acetamide;  
3,6-Bis-(morpholin-4-ylmethyl)-2,5-bis-(2-nitro-4-trifluoromethanesulfonyl-  
phenoxy)-benzene;  
[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-dimethyl-amine;  
*N*-(2-Ethylamino-5-trifluoromethanesulfonyl-phenyl)-2-(4-methanesulfonyl-phenyl)-  
acetamide;  
2-Hydroxy-5-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-terephthalic acid diethyl  
ester;  
{2-[(3-Trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetic acid;  
{3-[(3-Trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetic acid;  
{4-[(3-Trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetic acid;  
3,5-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-benzamide;  
3,5-Bis-(4-trifluoromethanesulfonyl-phenoxy)-benzoic acid;  
*N*-(4-Trifluoromethanesulfonyl-phenyl)-2-{2-[(4-trifluoromethanesulfonyl-  
phenylcarbamoyl)-methyl]-phenyl}-acetamide;  
*N*-(4-Trifluoromethanesulfonyl-phenyl)-2-{3-[(4-trifluoromethanesulfonyl-  
phenylcarbamoyl)-methyl]-phenyl}-acetamide;

*N*-(4-Trifluoromethanesulfonyl-phenyl)-2-{4-[(4-trifluoromethanesulfonyl-phenylcarbamoyl)-methyl]-phenyl}-acetamide;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-benzoic acid methyl ester;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-benzoic acid;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-*N*-pyridin-4-yl-benzamide;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-*N*-(4-methoxy-phenyl)-benzamide;

3-[4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-benzoylamino]-benzoic acid ethyl ester;

4-(1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-*N*-(2-pyrrolidin-1-yl-ethyl)-benzamide;

*N*-Ethyl-4-(1-ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazol-2-yl)-benzamide;

1-Ethyl-5-trifluoromethanesulfonyl-1*H*-benzoimidazole-2-carboxylic acid;

[2-(Benzoyl-butyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid methyl ester;

*N*-Benzyl-*N*-[butylcarbamoyl-(4-trifluoromethanesulfonyl-phenyl)-methyl]-benzamide;

*N*-[Butylcarbamoyl-(4-trifluoromethanesulfonyl-phenyl)-methyl]-*N*-(2-hydroxy-ethyl)-benzamide;

[2-(Acetyl-cyclopropyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

[2-(Acetyl-methyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

[2-(Benzoyl-cyclohexyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

*N*-Cyclohexyl-*N*-[(2,6-dimethyl-phenylcarbamoyl)-(4-trifluoromethanesulfonyl-phenyl)-methyl]-benzamide;

[2-(Acetyl-propyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

[2-(Acetyl-cyclohexyl-amino)-2-(4-trifluoromethanesulfonyl-phenyl)-acetylamino]-acetic acid ethyl ester;

{4-[4-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-benzenesulfonyl]-phenoxy}-acetic acid;

4-[2-(2-Nitro-4-trifluoromethanesulfonyl-phenoxy)-ethoxy]-benzoic acid;

2,5-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-terephthalic acid diethyl ester;

1-[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-piperidine;

4-[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-morpholine;

[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-(2-nitro-phenyl)-amine;

1-(2-Nitro-phenylamino)-3-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propan-2-ol;

[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propyl]-(4-nitro-phenyl)-amine;

1-(4-Nitro-phenylamino)-3-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propan-2-ol;

4-[2-Hydroxy-3-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propylamino]-benzenesulfonamide;

4-[2,3-Bis-(2-nitro-4-trifluoromethanesulfonyl-phenoxy)-propylamino]-benzenesulfonamide; or

Bis-{[4-(2-nitro-4-trifluoromethanesulfonyl)-phenoxy]-phenyl} sulfone.

44. (Original) The method of claim 39 wherein the compound is:

1,2-Bis(4-trifluoromethylsulfonamidophenyl)ethane;

1,2-Bis(2-methyl-4-trifluoromethylsulfonamidophenyl)ethane;

1,3-Bis(4-trifluoromethylsulfonamidophenoxy)-2,2-dimethylpropane;

1,3-Bis(4-trifluoromethylsulfonamidophenoxy)propane;

1,4-Bis(4-trifluoromethylsulfonamidophenoxy)butane;

1,4-Bis(4-trifluoromethylsulfonamidophenoxy)benzene;

1-(4-Aminophenoxy)-4-trifluoromethylsulfonamidophenoxy benzene;

Bis(4-trifluoromethylsulfonamidophenyl) ether;

1,3-Bis(4-trifluoromethylsulfonamidophenoxy)benzene;  
2,5-Bis(4-trifluoromethylsulfonamidophenyl)-(1,3,4)oxadiazole;  
Bis(4-trifluoromethylsulfonamidophenyl)-1,4-diisopropylbenzene;  
5-Trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid;  
1-Methyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid;  
(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetic acid;  
1-Methyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid phenylamide;  
5-Trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid phenylamide;  
3-[(1-Methyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carbonyl)-amino]-benzoic acid;  
3-[(5-Trifluoromethanesulfonylamino-1*H*-indole-2-carbonyl)-amino]-benzoic acid;  
4-[(5-Trifluoromethanesulfonylamino-1*H*-indole-2-carbonyl)-amino]-benzoic acid;  
4-[2-(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetylamino]-benzoic acid;  
3-[2-(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetylamino]-benzoic acid;  
4-{{[2-(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetylamino]-methyl}-benzoic acid;  
(2-Methyl-5-trifluoromethanesulfonylamino-indol-1-yl)-acetic acid *tert*-butyl ester;  
1-Methyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid ethyl ester;  
6-Trifluoromethanesulfonylamino-naphthalene-2-carboxylic acid;  
*N,N*-Bis[(6-carboxyl-naphthalen-2-yl)methyl] trifluoromethanesulfonamide;  
6-[(Methyl-trifluoromethanesulfonyl-amino)-methyl]-naphthalene-2-carboxylic acid;  
3-({6-[(Methyl-trifluoromethanesulfonyl-amino)-methyl]-naphthalene-2-carbonyl}-amino)-benzoic acid;  
1-*tert*-Butoxycarbonylmethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid ethyl ester;  
1-Carboxymethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid ethyl ester;  
1-*tert*-Butoxycarbonylmethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid;

1-Carboxymethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid;

1-Carboxymethyl-5-(*N,N*-ditrifluoromethanesulfonyl)amino-1*H*-indole-2-carboxylic acid ethyl ester;

1-*tert*-Butoxycarbonylmethyl-5-(*N,N*-ditrifluoromethanesulfonyl)amino-1*H*-indole-2-carboxylic acid ethyl ester;

1-Carboxymethyl-5-(*N,N*- ditrifluoromethanesulfonyl)amino-1*H*-indole-2-carboxylic acid;

1-Cyclohexylmethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid ethyl ester;

1-Benzyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid; or

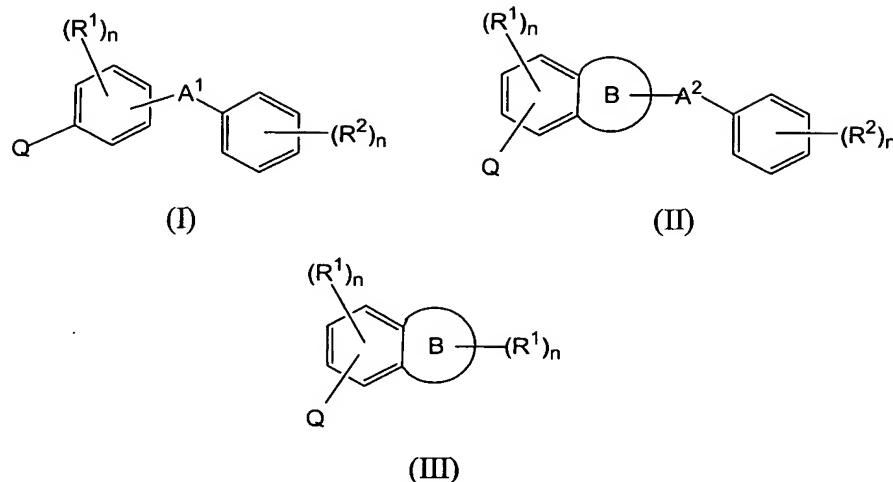
1-Cyclohexylmethyl-5-trifluoromethanesulfonylamino-1*H*-indole-2-carboxylic acid.

45. (Original) The method of claim 39 wherein said cancer is a solid tumor.

46. (Original) The method of claim 39 wherein said cancer is selected from the group consisting of glioma, melanoma, adenocarcinoma, Kaposi's sarcoma and hemangioma.

47. (Original) The method of claim 39 wherein said cancer is selected from the group consisting of ovarian, breast, lung, pancreatic, liver, prostate, colon, testicular, and epidermoid cancer.

48. (Original) A method for regulating, inhibiting or modulating protein tyrosine phosphatase signal transduction in a cell which comprises administering to the cell an effective amount of a compound having the formula:



or a pharmaceutically acceptable salt or solvate thereof, wherein:

$Q$  is  $CF_3SO_2$ ,  $CF_3SO_2NR^3$ ,  $CF_3SO_2R^4$  or  $CF_3SO_2N(R^3)R^4$ , wherein  $R^3$  is H, alkoxy, acyl or  $C_1$ - $C_3$  alkyl, each of which may be substituted or unsubstituted, and  $R^4$  is methylene which may be substituted or unsubstituted;

each  $R^1$  is independently  $C_1$ - $C_3$  alkyl,  $C_1$ - $C_3$  haloalkyl, CN,  $(C=O)OR$ ,  $(C=O)R^5$ , H, halo,  $NHR$ ,  $NH(C=O)OR$ ,  $NH(C=O)R^5$ ,  $NO_2$ ,  $NHSO_2R^5$ ,  $O(C=O)R$ , OR, OH,  $SO_2R^5$ ,  $R^4SO_2CF_3$  or tetrazole, wherein  $R^5$  is  $CF_3$ ,  $C_1$ - $C_3$  alkyl,  $NHR$  and wherein  $R$  is H,  $C_1$ - $C_3$  alkyl, aryl or heteroaryl, which may be substituted or unsubstituted;

each  $R^2$  is independently  $C_1$ - $C_3$  alkyl,  $C_1$ - $C_3$  haloalkyl, CN,  $(C=O)OR$ ,  $(C=O)R^5$ , H, halo,  $O(C=O)R$ , OR, OH,  $NHR$ ,  $NH(C=O)OR$ ,  $NH(C=O)R^5$ ,  $NO_2$ ,  $NHSO_2R^5$ ,  $SO_2R^5$ , tetrazole, or  $X_1$ - $R^6$ - $X_2$  wherein  $X_1$  may be present or absent and if present is O, N,  $(C=O)$ ,  $(C=O)NH$ ,  $NH(C=O)$ ,  $SO_2NH$ ,  $NHSO_2$ ,  $R^6$  is  $C_1$ - $3$  alkylene which may be substituted or unsubstituted and  $X_2$  is  $CF_3$ ,  $(C=O)OR$ ,  $(C=O)R^5$ , H,  $NH(C=O)R^5$ ,  $NH(C=O)OR$ ,  $NHSO_2R^5$ ,  $NRR^3$ ,  $O(C=O)R$ , OR,  $SO_2R^5$ , tetrazole;

each  $n$  is independently from 0 to 3;

ring  $B$  is an aryl, carbocyclic, heteroaryl, heterocyclic or phenyl ring which may be substituted or unsubstituted;

$A^1$  is a linkage in which the shortest path is 2-8 atoms in length wherein the atoms in the linkage are carbon which may be substituted or unsubstituted or the carbon replaced with a single nitrogen, oxygen or sulfur, or combination of nitrogen, oxygen and sulfur; the linkage may be or may contain an aryl, carbocyclic, heteroaryl, heterocyclic or a phenyl ring, which

may be directly in the linkage or appended to the linkage; the linkage may be acylalkyl, alkenylene, alkoxy, alkoxyalkyl, alkoxyamino, alkoxyarylalkoxy, alkoxyarylalkyl, alkoxyarylamino, alkoxyaryloxyalkyl, alkylamino, alkylaminoalkyl, alkylaminoarylaminolalkyl, alkylaryl, alkylarylalkyl, alkylarylamino, alkylaryloxy, alkylene, alkylenediamine, alkylenedioxy, alkyloxy, alkyloxyaryl, alkyloxyarylalkyloxy, alkyloxyaryloxyalkyl, alkylsulfonylamino, alkylthio, alkylthioalkyl, alkynylene, N-sulfonamido, N-amido, aminoalkyl, aminoalkylamino, aminoalkylarylalkyl, aminoalkylarylalkylamino, aminoalkylaryloxy, aminoalkyloxy, aminoaryl, aminoarylalkyl, aminoarylcarbonyl, aminoaryloxy, aminoaryloxyalkyl, aminoarylsulfonyl, aryl, arylamino, aryldioxy, aryldiamine, aryloxy, aryloxyalkyl, aryloxyamino, aryloxyaminoalkyl, aryloxycarbonyl, aryloxysulfonyl, benzimidazole, benzo[b]furan, benzo[b]thiophene, C-amido, carbonylarylaminol, carbonylarylcarbonyl, carbonylaryloxy, chromene, cycloalkylene, disulfide, furan, haloalkyl, imidazole, imidazolidine, imidazoline, indole, isothiazole, isoxazole, morpholine, oxadiazole, oxazole, oxirane, parathiazine, phenothiazine, piperazine, piperidine, purine, pyran, pyrazine, pyrazole, pyrazolidine, pyrimidine, pyridine, pyrrole, pyrrolidine, quinoline, sulfonamido, sulfonylalkyl, sulfonylarylaminol, sulfonylaryloxy, sulfonylarylsulfonyl, thiadiazole, thiazole, thiophene, triazine, triazole, unsubstituted azeridine, ureido, which may be substituted or unsubstituted;

$A^2$  is a linkage in which the shortest path is 0-6 atoms in length wherein the atoms in the linkage are carbon which may be substituted or unsubstituted or the carbon replaced with a single nitrogen, oxygen or sulfur, or combination of nitrogen, oxygen and sulfur; the linkage may be or may contain an aryl, carbocyclic, heteroaryl, heterocyclic or a phenyl ring, which may be directly in the linkage or appended to the linkage; the linkage may be single atom C, O, S or N which may be substituted or unsubstituted; the linkage may be acylalkyl, alkenylene, alkoxy, alkoxyalkyl, alkoxyamino, alkoxyarylalkoxy, alkoxyarylalkyl, alkoxyarylamino, alkoxyaryloxyalkyl, alkylamino, alkylaminoalkyl, alkylaminoarylaminolalkyl, alkylaryl, alkylarylalkyl, alkylarylamino, alkylaryloxy, alkylene, alkylenediamine, alkylenedioxy, alkyloxy, alkyloxyaryl, alkyloxyarylalkyloxy, alkyloxyaryloxyalkyl, alkylsulfonylamino, alkylthio, alkylthioalkyl, alkynylene, N-sulfonamido, N-amido, aminoalkyl, aminoalkylamino, aminoalkylarylalkyl, aminoalkylarylalkylamino, aminoalkylaryloxy, aminoalkyloxy, aminoaryl, aminoarylalkyl,

aminoarylcarbonyl, aminoaryloxy, aminoaryloxyalkyl, aminoarylsulfonyl, aryl, arylamino, ortho or para aryldioxy, substituted meta-aryldioxy, aryldiamine, aryloxy, aryloxyalkyl, aryloxyamino, aryloxyaminoalkyl, aryloxycarbonyl, aryloxysulfonyl, benzimidazole, benzo[b]furan, benzo[b]thiophene, C-amido, carbonylarylarnino, carbonylarylcarbonyl, carbonylaryloxy, chromene, cycloalkylene, furan, haloalkyl, imidazole, imidazolidine, imidazoline, indole, isothiazole, isoxazole, morpholine, oxadiazole, oxazole, oxirane, parathiazine, phenothiazine, piperazine, piperidine, purine, pyran, pyrazine, pyrazole, pyrazolidine, pyrimidine, pyridine, pyrrole, pyrrolidine, quinoline, sulfonamido, sulfonylalkyl, sulfonylarylarnino, sulfonylaryloxy, sulfonylarylsulfonyl, thiadiazole, thiazole, thiophene, triazine, triazole, unsubstituted azeridine, ureido, which may be substituted or unsubstituted

wherein said compound regulates, inhibits or modulates protein tyrosine phosphatase signal transduction.

49 – 89. (Canceled)